



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:25 pm BST

PDB ID : 6JY1
Title : Crystal Structure of a Group II pyridoxal dependent decarboxylase, LLP-bound form from Methanocaldococcus jannaschii at 1.72 Å
Authors : Manoj, N.; Gayathri, S.C.
Deposited on : 2019-04-25
Resolution : 1.72 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

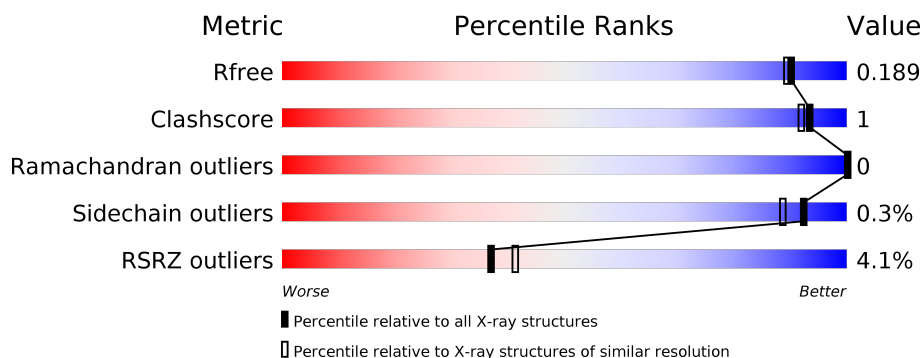
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.72 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5722 (1.74-1.70)
Clashscore	141614	6152 (1.74-1.70)
Ramachandran outliers	138981	6051 (1.74-1.70)
Sidechain outliers	138945	6051 (1.74-1.70)
RSRZ outliers	127900	5629 (1.74-1.70)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	415	<div> <div>4%</div> <div> <div></div> <div>90%</div> <div>• 6%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6591 atoms, of which 3147 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called L-tyrosine/L-aspartate decarboxylase.

Mol	Chain	Residues	Atoms								ZeroOcc	AltConf	Trace
1	A	390	Total	C	H	N	O	P	S		0	6	0
			6257	2004	3129	520	587	1	16				

There are 19 discrepancies between the modelled and reference sequences:

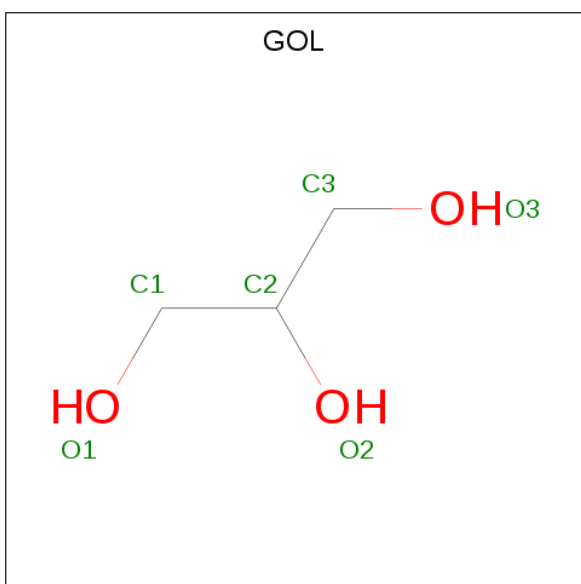
Chain	Residue	Modelled	Actual	Comment	Reference
A	-18	MET	-	initiating methionine	UNP Q60358
A	-17	GLY	-	expression tag	UNP Q60358
A	-16	SER	-	expression tag	UNP Q60358
A	-15	ASP	-	expression tag	UNP Q60358
A	-14	LYS	-	expression tag	UNP Q60358
A	-13	ILE	-	expression tag	UNP Q60358
A	-12	HIS	-	expression tag	UNP Q60358
A	-11	HIS	-	expression tag	UNP Q60358
A	-10	HIS	-	expression tag	UNP Q60358
A	-9	HIS	-	expression tag	UNP Q60358
A	-8	HIS	-	expression tag	UNP Q60358
A	-7	HIS	-	expression tag	UNP Q60358
A	-6	GLU	-	expression tag	UNP Q60358
A	-5	ASN	-	expression tag	UNP Q60358
A	-4	LEU	-	expression tag	UNP Q60358
A	-3	TYR	-	expression tag	UNP Q60358
A	-2	PHE	-	expression tag	UNP Q60358
A	-1	GLN	-	expression tag	UNP Q60358
A	0	GLY	-	expression tag	UNP Q60358

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

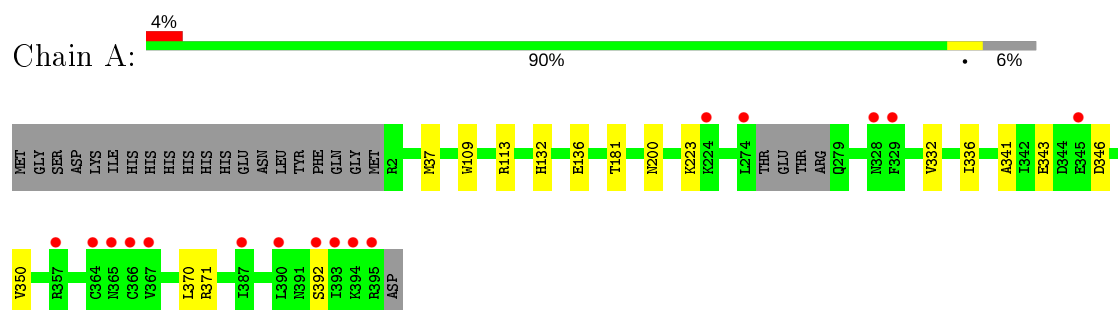


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	H	O	0	0
			14	3	8	3		
3	A	1	Total	C	H	O	0	0
			11	3	5	3		
3	A	1	Total	C	H	O	0	0
			11	3	5	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	273	Total	O	0	0
			273	273		

- Molecule 1: L-tyrosine/L-aspartate decarboxylase



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	98.58 Å 98.58 Å 122.28 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.72 – 1.72 51.96 – 1.72	Depositor EDS
% Data completeness (in resolution range)	99.9 (45.72-1.72) 99.9 (51.96-1.72)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 1.72 Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.163 , 0.188 0.163 , 0.189	Depositor DCC
R_{free} test set	3191 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.44 , 49.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	6591	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.26% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, LLP, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.44	0/3163	0.62	0/4262

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3128	3129	3142	9	0
2	A	25	0	0	0	0
3	A	18	18	24	1	0
4	A	273	0	0	1	0
All	All	3444	3147	3166	9	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 1.

All (9) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:223:LYS:HD3	1:A:336:ILE:HD13	1.96	0.47
1:A:132:HIS:HD2	1:A:181[B]:THR:HG21	1.80	0.47
1:A:332:VAL:HG22	1:A:343:GLU:HG3	1.97	0.47
1:A:37[A]:MET:HE2	1:A:371:ARG:HH12	1.81	0.45
1:A:346:ASP:O	1:A:350:VAL:HG23	2.16	0.45
1:A:136:GLU:HG2	4:A:561:HOH:O	2.19	0.42
1:A:341:ALA:HA	1:A:370:LEU:O	2.20	0.42
1:A:109:TRP:CZ2	1:A:113:ARG:HG3	2.54	0.42
1:A:200:ASN:C	3:A:408:GOL:H11	2.42	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	391/415 (94%)	376 (96%)	15 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	335/363 (92%)	334 (100%)	1 (0%)	92	89

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	392	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

1 non-standard protein/DNA/RNA residue is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	LLP	A	245	1	23,24,25	2.28	5 (21%)	25,32,34	1.35	3 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	LLP	A	245	1	-	2/16/17/19	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	245	LLP	C4-C4'	6.55	1.59	1.46
1	A	245	LLP	C4'-NZ	4.52	1.42	1.27
1	A	245	LLP	C4-C5	-4.47	1.36	1.42
1	A	245	LLP	CB-CA	-2.57	1.50	1.53
1	A	245	LLP	C2'-C2	2.06	1.53	1.50

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	245	LLP	C4-C4'-NZ	-3.53	108.10	124.31
1	A	245	LLP	CE-NZ-C4'	-2.59	110.95	118.90
1	A	245	LLP	CG-CD-CE	-2.04	106.46	113.57

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	245	LLP	C4-C4'-NZ-CE
1	A	245	LLP	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	402	-	4,4,4	0.16	0	6,6,6	0.09	0
3	GOL	A	408	-	5,5,5	1.02	0	5,5,5	0.96	0
2	SO4	A	403	-	4,4,4	0.18	0	6,6,6	0.14	0
2	SO4	A	404	-	4,4,4	0.12	0	6,6,6	0.16	0
3	GOL	A	406	-	5,5,5	0.91	0	5,5,5	0.94	0
2	SO4	A	405	-	4,4,4	0.17	0	6,6,6	0.44	0
2	SO4	A	401	-	4,4,4	0.16	0	6,6,6	0.18	0
3	GOL	A	407	-	5,5,5	1.13	0	5,5,5	0.68	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	406	-	-	0/4/4/4	-
3	GOL	A	408	-	-	2/4/4/4	-
3	GOL	A	407	-	-	4/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	408	GOL	O1-C1-C2-C3
3	A	407	GOL	C1-C2-C3-O3
3	A	407	GOL	O1-C1-C2-C3
3	A	407	GOL	O1-C1-C2-O2
3	A	408	GOL	O1-C1-C2-O2
3	A	407	GOL	O2-C2-C3-O3

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	408	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	389/415 (93%)	0.19	16 (4%) 37 41	15, 25, 55, 89	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	366	CYS	10.7
1	A	393	ILE	4.8
1	A	395	ARG	4.2
1	A	367	VAL	3.7
1	A	224	LYS	3.3
1	A	357	ARG	3.3
1	A	394	LYS	3.0
1	A	390	LEU	2.9
1	A	274	LEU	2.9
1	A	345	GLU	2.7
1	A	328	ASN	2.7
1	A	365	ASN	2.6
1	A	387	ILE	2.5
1	A	392	SER	2.2
1	A	364	CYS	2.2
1	A	329	PHE	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	LLP	A	245	24/25	0.98	0.11	16,23,35,36	0

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	A	406	6/6	0.78	0.18	40,50,64,69	0
3	GOL	A	408	6/6	0.83	0.27	26,35,47,47	11
2	SO4	A	404	5/5	0.87	0.26	52,67,91,106	0
2	SO4	A	405	5/5	0.88	0.13	41,54,59,59	0
2	SO4	A	403	5/5	0.92	0.33	55,68,76,90	0
3	GOL	A	407	6/6	0.94	0.21	27,46,63,63	0
2	SO4	A	402	5/5	0.95	0.23	37,71,87,99	0
2	SO4	A	401	5/5	0.98	0.10	50,53,65,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.